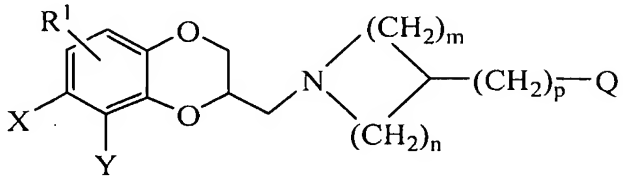
This listing of claims will replace all prior versions, and listings, of claims in the application.

Listing of Claims

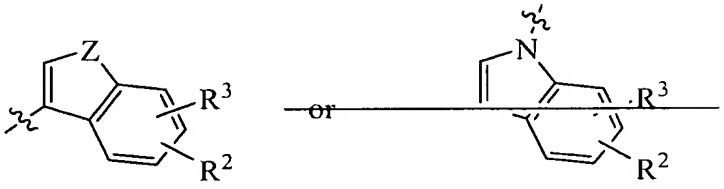
1. ***(currently amended)*** A compound of Formula I:



I

wherein

Q is



a **b**

R^1, R^2 and R^3 are, independently, hydrogen, hydroxy, halo, cyano, carboxamido, carboalkoxy of two to six carbon atoms, trifluoromethyl, alkyl of 1 to 6 carbon atoms, alkoxy of 1 to 6 carbon atoms, alkanoyl of 2 to 6 carbon atoms, alkanoyloxy of 2 to 6 carbon atoms, amino, mono- or di-alkylamino in which each alkyl group has 1 to 6 carbon atoms, alkanamido of 2 to 6 carbon atoms, alkanesulfonyl of 1 to 6 carbon atoms or alkanesulfonamido of 1 to 6 carbon atoms;

X and Y ~~are, independently, hydrogen, hydroxy, halo, cyano, carboxamido, carboalkoxy of two to six carbon atoms, trifluoromethyl, alkyl of 1 to 6 carbon atoms, alkoxy of 1 to 6 carbon atoms, alkanoyl of 2 to 6 carbon atoms, alkanoyloxy of 2 to 6 carbon atoms, amino, mono- or di-alkylamino in which each alkyl group has 1 to 6 carbon atoms, alkanamido of 2 to 6 carbon~~

~~**atoms, alkanesulfonyl of 1 to 6 carbon atoms or alkanesulfonamido of 1 to 6 carbon atoms, or X and Y**~~, taken together, form ~~**-N=C(R^4)-C(R^5)=N-,**~~ -N=C(R^4)-C(R^6)=CH-, ~~**-N=C(R^4)-N=CH-, -N=C(R^4)-O-, -NH-C(R^7)=N- or -NH-C(R^8)=CH-**~~;

R^4 and R^5 are, independently, hydrogen, halo, amino, mono- or di-alkylamino in which each alkyl group has 1 to 6 carbon atoms or alkyl of 1 to 6 carbon atoms;

R^6 is hydrogen or alkyl of 1 to 6 carbon atoms;

~~**R^7 is hydrogen, halo, trifluoromethyl, pentafluoroethyl, amino, mono- or di-alkylamino in which each alkyl group has 1 to 6 carbon atoms or alkyl of 1 to 6 carbon atoms;**~~

~~**R^8 is hydrogen, halo, trifluoromethyl, pentafluoroethyl, or alkyl of 1 to 6 carbon atoms;**~~

Z is O, S, or NR^9, in which R^9 is hydrogen or alkyl of 1 to 6 carbon atoms;

n is an integer ~~**0,**~~ 1~~,~~ or 2;

m is an integer from 1 to ~~**4**~~ **2, provided that m + n ≤ 4 and that when m = n = 2, and Q is b then X and Y are not NH-C(R^8)=CH-**; and

p is an integer from 1 to ~~**3**~~ **2**, provided that p + n is 2 or 3;

or a pharmaceutically acceptable salt thereof.

2-3. ***(cancelled)***

4. ***(original)*** A compound according to claim 1, wherein Z is NR^9 or a pharmaceutically acceptable salt thereof.

5. ***(currently amended)*** A compound according to claim 1, wherein n is ~~**0 or**~~ 1 or a pharmaceutically acceptable salt thereof.

6. ***(currently amended)*** A compound according to claim 1, wherein m is 1 ~~**to 3**~~ or a pharmaceutically acceptable salt thereof.

7. ***(currently amended)*** A compound according to claim 1, wherein p is 1 ~~**or 2**~~ or a pharmaceutically acceptable salt thereof.

8. ***(original)*** A compound according to claim 1, wherein R^1 is hydrogen, halo, cyano, trifluoromethyl, alkyl of 1 to 6 carbon atoms or alkoxy of 1 to 6 carbon atoms or a pharmaceutically acceptable salt thereof.

9. ***(original)*** A compound according to claim 1, wherein R^2 and R^3 are independently selected from hydrogen, hydroxy, halo, cyano, carboxamido, alkyl of 1 to 6 carbon atoms, or alkoxy of 1 to 6 carbon atoms or a pharmaceutically acceptable salt thereof.

10. ***(original)*** A compound according to claim 1, wherein R^4 and R^5 are independently hydrogen, amino or alkyl of 1 to 6 carbon atoms or a pharmaceutically acceptable salt thereof.

11. ***(cancelled)***

12. ***(currently amended)*** A compound according to claim 1, wherein R^6 is hydrogen or alkyl of 1 to 3 carbon atoms, Z is NR^9 in which R^9 is hydrogen or alkyl of 1 to 3 carbon atoms, ~~**n is 0 or 1, m is 1 to 3 and p is 1 or 2**~~ or a pharmaceutically acceptable salt thereof.

13. ***(cancelled)***

14. ***(original)*** A compound according to claim 1, wherein said compound is 2-[3-(5-fluoro-1H-indol-3-ylmethyl)-azetidin-1-ylmethyl]-8-methyl-2,3-dihydro-[1,4]dioxino[2,3-f]quinoline or a pharmaceutically acceptable salt thereof.

15-16. ***(cancelled)***

17. ***(original)*** A compound according to claim 1, wherein said compound is 2-[3-(1H-indol-3-ylmethyl)-azetidin-1-ylmethyl]-8-methyl-2,3-dihydro-[1,4]dioxino[2,3-f]quinoline or a pharmaceutically acceptable salt thereof.

18. ***(original)*** A compound according to claim 1, wherein said compound is 2-[3-(5-fluoro-1-methyl-1H-indol-3-ylmethyl)-azetidin-1-ylmethyl]-8-methyl-2,3-dihydro-[1,4]dioxino[2,3-f]quinoline or a pharmaceutically acceptable salt thereof.

19-24. ***(cancelled)***

25. ***(original)*** A compound according to claim 1, wherein said compound is 8-Methyl-2-[3-(5-methyl-1H-indol-3-ylmethyl)-azetidin-1-ylmethyl]-2,3-dihydro-[1,4]dioxino[2,3-f]quinoline or a pharmaceutically acceptable salt thereof.

26. ***(original)*** A compound according to claim 1, wherein said compound is the S enantiomer at the 2-aminomethyl-2,3-dihydro-1,4-benzodioxan moiety, substantially free of the R enantiomer of said compound.

27-29. ***(cancelled)***

30. ***(original)*** A pharmaceutical composition, comprising:
an effective amount of a compound according to claim 1; and
a pharmaceutically acceptable carrier or excipient.

31. ***(new)*** A compound selected from the group consisting of:
2-[3-(5-fluoro-1H-indol-3-ylmethyl)-piperidin-1-ylmethyl]-8-methyl-2,3-dihydro-[1,4]dioxino[2,3-f]quinoline;
2-[3-(6-fluoro-1H-indol-3-ylmethyl)-piperidin-1-ylmethyl]-8-methyl-2,3-dihydro-[1,4]dioxino[2,3-f]quinoline;

2-({4-[(6-fluoro-1H-indol-1-yl)methyl]piperidin-1-yl})-8-methyl-2,3-dihydro[1,4]dioxino[2,3-f]quinoline;

2-({4-[(6-fluoro-1H-indol-1-yl)ethyl]piperidin-1-yl})-8-ethyl-2,3-dihydro[1,4]dioxino[2,3-f]quinoline;

1-[(1-{[8-methyl-2,3-dihydro[1,4]-dioxino[2,3-f]quinolin-2-yl]methyl})piperidin-4-yl]-1H-indole-6-carbonitrile;

2-[3-(6-fluoro-indol-1-ylmethyl)-azetidin-1-ylmethyl]-8-methyl-2,3-dihydro-[1,4]dioxino[2,3-*f*]quinoline;

2-{3-[2-(6-fluoro-indol-1-yl)-ethyl]-azetidin-1-ylmethyl}-8-methyl-2,3-dihydro-[1,4]dioxino[2,3-f]quinoline;

1-{2-[1-(8-methyl-2,3-dihydro-[1,4]-dioxino[2,3-f]quinolin-2-ylmethyl)-azetinin-3-yl]-ethyl}-1H-indole-6-carbonitrile; and

pharmaceutically acceptable salts thereof.